

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSPTANXR1625

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

FILE 'HOME' ENTERED AT 11:16:18 ON 02 JUL 2007

FILE 'REGISTRY' ENTERED AT 11:16:27 ON 02 JUL 2007
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2007 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 1 JUL 2007 HIGHEST RN 940612-32-8
DICTIONARY FILE UPDATES: 1 JUL 2007 HIGHEST RN 940612-32-8

New CAS Information Use Policies, enter HELP USAGETERMS for details.

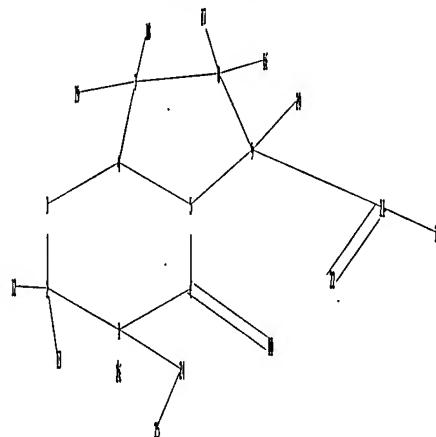
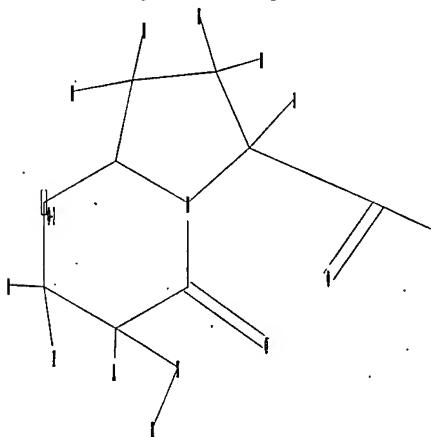
TSCA INFORMATION NOW CURRENT THROUGH December 2, 2006

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stnqgen/stndoc/properties.html>

=>
Uploading C:\Program Files\Stnexp\Queries\10563288a.str



chain nodes :

10 12 13 16 17 18 19 20 21 22 23 24 25 26

ring nodes :

1 2 3 4 5 6 7 8 9

chain bonds :
1-24 1-26 2-12 2-13 6-10 7-18 7-19 8-16 8-17 9-20 9-21 21-22 21-23

24-25

ring bonds :

1-2 1-6 2-3 3-4

exact/norm bonds :

1-2 1-6 1-2

exact bonds :

2-12 2-13 4-7 7-8 7

containing 1 :

Match level :

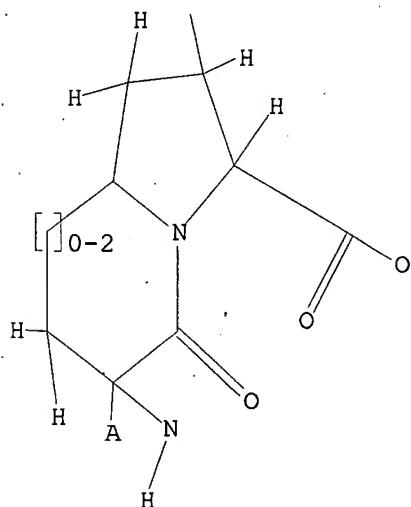
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
12:CLASS 13:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS 20:CLASS 21:CLASS
22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS

L1 STRUCTURE UPLOADED

=> d 11

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 11 full

FULL SEARCH INITIATED 11:16:46 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 2321 TO ITERATE

100.0% PROCESSED 2321 ITERATIONS
SEARCH TIME: 00.00.01

43 ANSWERS

L2 43 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE

ENTRY

TOTAL

FULL ESTIMATED COST

SESSION

172.10

172.31

FILE 'CAPLUS' ENTERED AT 11:16:54 ON 02 JUL 2007

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2007 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the

American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 2 Jul 2007 VOL 147 ISS 2
FILE LAST UPDATED: 1 Jul 2007 (20070701/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply.
They are available for your review at:

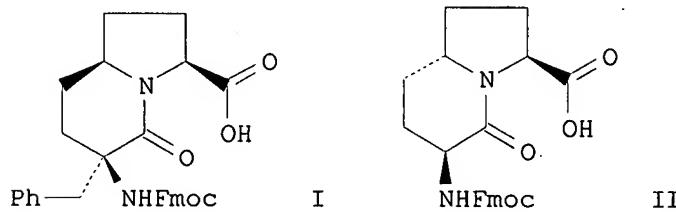
<http://www.cas.org/infopolicy.html>

=> s 12 full
L3 18 L2

=> s 13 and py<2003
22885672 PY<2003
L4 13 L3 AND PY<2003

=> d ibib abs hitstr tot

L4 ANSWER 1 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN
ACCESSION NUMBER: 2003:69868 CAPLUS
DOCUMENT NUMBER: 139:53275
TITLE: Cyclic RGD peptides containing azabicycloalkane reverse-turn mimics
AUTHOR(S): Belvisi, Laura; Caporale, Andrea; Colombo, Matteo; Manzoni, Leonardo; Potenza, Donatella; Scolastico, Carlo; Castorina, Massimo; Cati, Matilde; Giannini, Giuseppe; Pisano, Claudio
CORPORATE SOURCE: Dipartimento di Chimica Organica e Industriale, Universita degli Studi di Milano, Milan, I-20133, Italy
SOURCE: Helvetica Chimica Acta (2002), 85(12), 4353-4368
CODEN: HCACAV; ISSN: 0018-019X
PUBLISHER: Verlag Helvetica Chimica Acta
DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 139:53275
GI



AB The Fmoc-protected lactams I and II were used to prepare cyclo(Arg-Gly-Asp-lactam) (III) and cyclo(Arg-Gly-Asp-Phe-lactam), which contain the Arg-Gly-Asp (RGD) recognition motif. Their solid-phase synthesis, conformational anal., and binding to purified $\alpha v\beta 3$ and $\alpha v\beta 5$ integrins are reported. Compound III was found to act as an active and selective inhibitor of the $\alpha v\beta 5$ integrin.

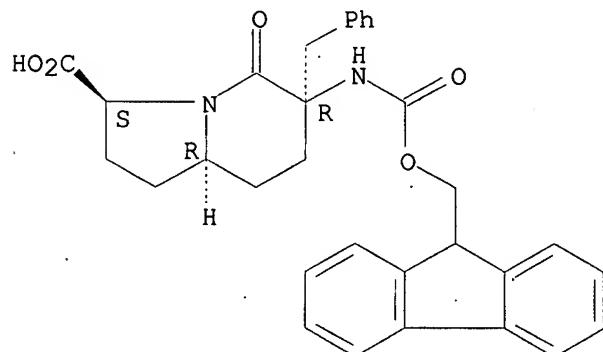
IT 220719-80-2

RL: RCT (Reactant); RACT (Reactant or reagent)
(solid-phase synthesis, conformational anal., and integrin-binding of cyclic RGD peptides containing azabicycloalkane reverse-turn mimics)

RN 220719-80-2 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[[[9H-fluoren-9-ylmethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, (3S,6R,8aR)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



REFERENCE COUNT: 49 THERE ARE 49 CITED REFERENCES AVAILABLE FOR THIS RECORD: ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 2 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1999:529170 CAPLUS

DOCUMENT NUMBER: 131:157992

TITLE: Synthesis of β -sheet mimetics as inhibitors of protease, kinase, transcription factors, and protein-protein binding interactions

INVENTOR(S): Gaber, Maher N.; McMillian, Michael K.; Kahn, Michael S.; Tulinsky, John E.; Mathew, Jessymol

PATENT ASSIGNEE(S): Molecumetics Ltd., USA

SOURCE: PCT Int. Appl., 279 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

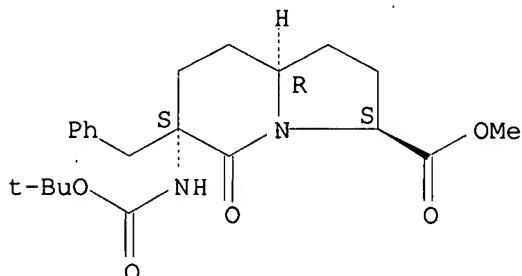
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9941276	A1	19990819	WO 1998-US2891	19980212 <--
W: AL, AM, AT, AU, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, GM, GW, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
CA 2319766	A1	19990819	CA 1998-2319766	19980212 <--
AU 9866557	A	19990830	AU 1998-66557	19980212 <--
AU 748887	B2	20020613		
EP 1053246	A1	20001122	EP 1998-908551	19980212 <--
EP 1053246	B1	20030102		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, FI				
JP 2002503674	T	20020205	JP 2000-531467	19980212 <--
AT 230414	T	20030115	AT 1998-908551	19980212
NZ 505980	A	20030131	NZ 1998-505980	19980212
ES 2192764	T3	20031016	ES 1998-908551	19980212
PRIORITY APPLN. INFO.:			WO 1998-US2891	A 19980212
OTHER SOURCE(S):	MARPAT	131:157992		
GI				

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

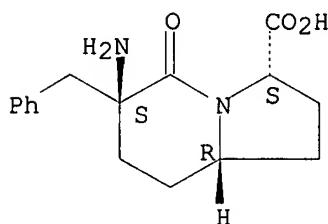
- AB β -Sheet mimetics I [A = CO, (CH₂)₀₋₄, CO(CH₂)₁₋₃, (CH₂)₁₋₂₀, (CH₂)_{1-2S}; B = N, CH; C = CO, CO(CH₂)₁₋₃, (CH₂)₀₋₃, O, S, O(CH₂)₁₋₂, S(CH₂)₁₋₂; D = N, CR₄; E = CR₁(NH₂), NZ, CZR₁; F is an optional carbonyl moiety; R₁, R₃, R₄ is an amino acid side chain moiety or derivative; R₂, R_{2'} are ring substituents selected from an amino acid side chain moiety or derivative, where at least one R₂ together with C forms a fused (un)substituted homo- or heterocyclic ring; Y, Z represent the remainder of the mol.; any adjacent CH groups of the bicyclic ring may form a double bond] were prepared as inhibitors of protease, kinase, transcription factors, and protein-protein binding interactions. Thus, arginine derivative II was prepared and assayed for inhibition of various enzymes (IC₅₀ = 1.2 and 140 nM for thrombin and Factor VII, resp.).
- IT 183442-92-4P 183442-93-5P 183624-03-5P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(synthesis of β -sheet mimetics as inhibitors of protease, kinase, transcription factors, and protein-protein binding interactions)
- RN 183442-92-4 CAPLUS
- CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, methyl ester, (3R,6R,8aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



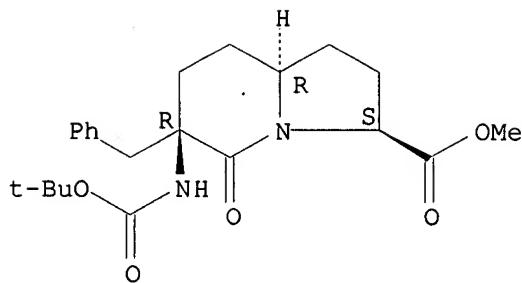
- RN 183442-93-5 CAPLUS
CN 3-Indolizinecarboxylic acid, 6-aminoctahydro-5-oxo-6-(phenylmethyl)-, (3R,6R,8aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



- RN 183624-03-5 CAPLUS
CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, methyl ester, (3R,6S,8aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 3 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1999:396495 CAPLUS

DOCUMENT NUMBER: 131:200003

TITLE: Synthesis of dipeptide secondary structure mimetics

AUTHOR(S): Eguchi, Masakatsu; Kim, Hwa-Ok; Gardner, Benjamin S.; Boatman, P. Douglas; Lee, Min S.; Nakanishi, Hiroshi; Kahn, Michael

CORPORATE SOURCE: Molecumetics Ltd., Bellevue, WA, 98005, USA

SOURCE: Peptides: Frontiers of Peptide Science, Proceedings of the American Peptide Symposium, 15th, Nashville, June 14-19, 1997 (1999), Meeting Date 1997, 212-213. Editor(s): Tam, James P.; Kaumaya, Pravin T. P. Kluwer: Dordrecht, Neth.

CODEN: 67UCAR

DOCUMENT TYPE: Conference

LANGUAGE: English

AB A symposium with four refs. on the synthesis and biol. activity of two diastereomeric title compds.

IT 203453-44-5P 203455-60-1P

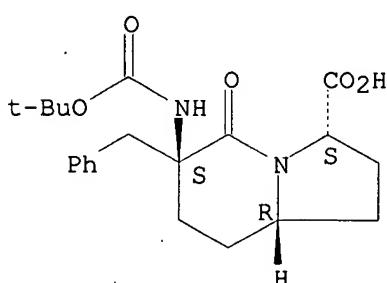
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation and reaction of in the synthesis of dipeptide secondary structure mimetics as thrombin inhibitors)

RN 203453-44-5 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, (3S,6S,8aR)- (9CI) (CA INDEX NAME)

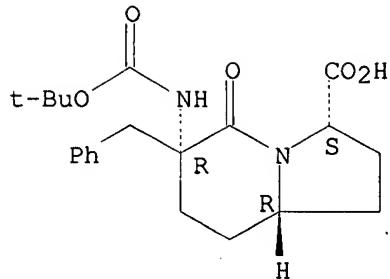
Absolute stereochemistry.



RN 203455-60-1 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, (3S,6R,8aR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 4 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1999:89647 CAPLUS

DOCUMENT NUMBER: 130:182766

TITLE: Conformational preferences of peptides containing reverse-turn mimetic bicyclic lactams. Inverse γ -turns versus type-II' β -turns. Insights into β -hairpin stability

AUTHOR(S): Belvisi, Laura; Gennari, Cesare; Mielgo, Antonia; Potenza, Donatella; Scolastico, Carlo

CORPORATE SOURCE: Dipartimento Chimica Organica Industriale, Univ. Studi Milano, Milan, Italy

SOURCE: European Journal of Organic Chemistry. (1999), (2), 389-400

CODEN: EJOCFK; ISSN: 1434-193X

PUBLISHER: Wiley-VCH Verlag GmbH

DOCUMENT TYPE: Journal

LANGUAGE: English

AB The conformational preferences of constrained peptides containing reverse-turn mimetic bicyclic lactams were investigated by NMR and IR. The exptl. results were complemented by computer modeling studies and show that the constrained peptides form an inverse γ -turn or a type-II' β -turn through intramol. H-bonding, depending on the nature of the reverse-turn mimic. In N-acetylated tetrapeptide mimics incorporating the two different bicyclic lactams, H(5) is available for either a γ -turn (7-membered ring with the CO group of the bicyclic lactam) or a β -turn (10-membered ring with the CO group of residue 2). Peptides incorporating a (5,7)-bicyclic lactam predominantly induce the γ -turn conformation, while those incorporating a (5,6)-bicyclic lactam can promote either a γ -turn or a β -turn conformation, with the β -turn usually being preferred and with varying degrees of β -hairpin formation.

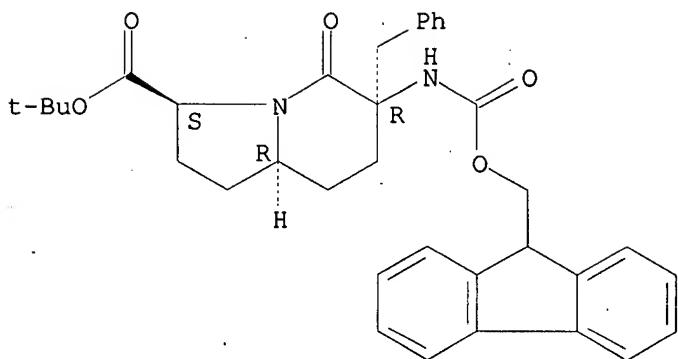
IT 220563-70-2

RL: PRP (Properties)
(conformational anal.)

RN 220563-70-2 CAPLUS

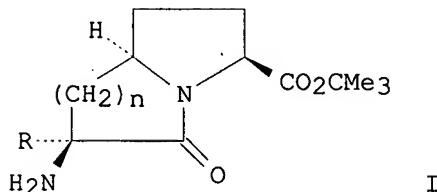
CN 3-Indolizinecarboxylic acid, 6-[[[(9H-fluoren-9-ylmethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, 1,1-dimethylethyl ester, (3S,6R,8aR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).



REFERENCE COUNT: 68 THERE ARE 68 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 5 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN
 ACCESSION NUMBER: 1999:89646 CAPLUS
 DOCUMENT NUMBER: 130:196946
 TITLE: Solid-phase synthesis of peptides containing reverse-turn mimetic bicyclic lactams
 AUTHOR(S): Gennari, Cesare; Mielgo, Antonia; Potenza, Donatella; Scolastico, Carlo; Piarulli, Umberto; Manzoni, Leonardo
 CORPORATE SOURCE: Dipartimento Chimica Organica Industriale, Univ. Studi Milano, Milan, Italy
 SOURCE: European Journal of Organic Chemistry (1999), (2), 379-388
 CODEN: EJOCFK; ISSN: 1434-193X
 PUBLISHER: Wiley-VCH Verlag GmbH
 DOCUMENT TYPE: Journal
 LANGUAGE: English
 GI



I

AB The solid-phase synthesis and characterization of a series of peptides containing reverse-turn mimetic bicyclic lactams I ($n = 3$, $R = H$; $n = 2$, $R = PhCH_2$) is reported. The bicyclic lactams possess high structural similarity to the 2 central residues of a β -turn. Amino acid conjugates of these bicyclic lactams were synthesized on solid supports following a 9-fluorenylmethoxycarbonyl (FMOC) protection strategy on Wang-Merrifield resin. Coupling between amino acids was accomplished by diisopropylcarbodiimide (DIC)/hydroxyazabenzotriazole (HOAt). Coupling between amino acids and the mimics was performed with the potent Carpino's reagent, O-(7-azabenzotriazol-1-yl)-N,N,N',N'-tetramethyluronium hexafluorophosphate (HATU). The final compds. were cleaved from the resin and obtained as N-acetylated Me esters or benzyl amides.

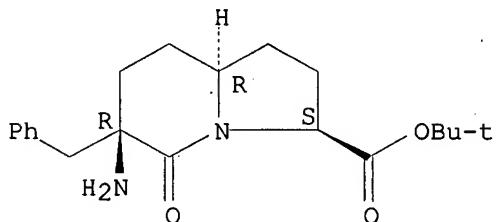
IT 208455-27-0

RL: RCT (Reactant); RACT (Reactant or reagent)
 (solid-phase synthesis of peptides containing reverse-turn mimetic bicyclic lactams)

RN 208455-27-0 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-aminoctahydro-5-oxo-6-(phenylmethyl)-, 1,1-dimethylethyl ester, (3S,6R,8aR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



IT 220563-70-2P 220719-80-2P

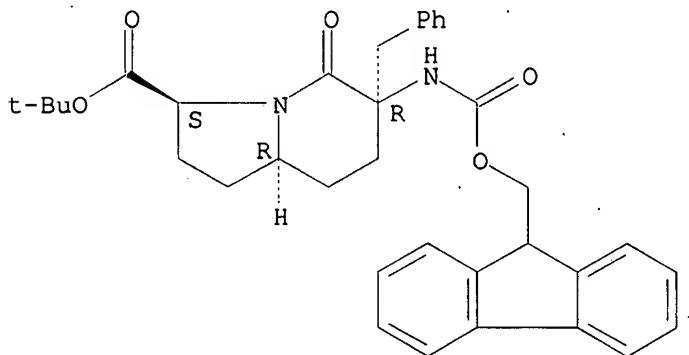
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(solid-phase synthesis of peptides containing reverse-turn mimetic bicyclic lactams)

RN 220563-70-2 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[(9H-fluoren-9-ylmethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, 1,1-dimethylethyl ester, (3S,6R,8aR)- (9CI) (CA INDEX NAME)

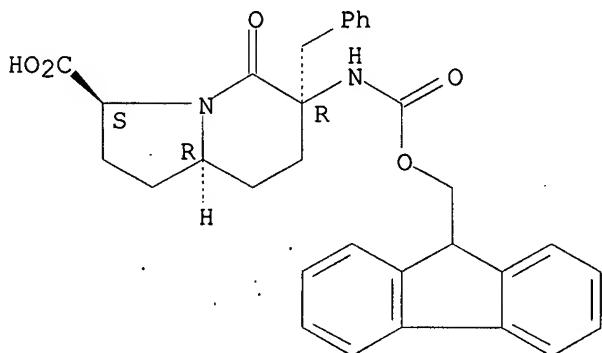
Absolute stereochemistry. Rotation (+).



RN 220719-80-2 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[(9H-fluoren-9-ylmethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, (3S,6R,8aR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



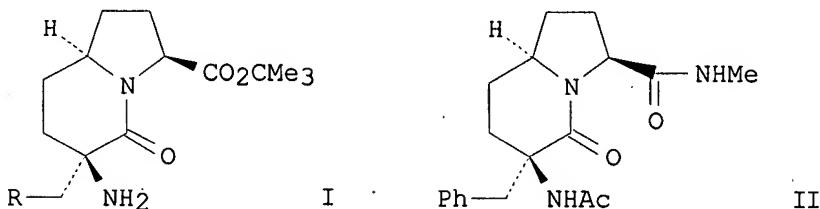
REFERENCE COUNT:

90

THERE ARE 90 CITED REFERENCES AVAILABLE FOR THIS

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 6 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN
 ACCESSION NUMBER: 1998:304016 CAPLUS
 DOCUMENT NUMBER: 129:41402
 TITLE: Stereoselective synthesis of 6,5-bicyclic reverse-turn peptidomimetics
 AUTHOR(S): Colombo, Lino; Di Giacomo, Marcello; Brusotti, Gloria;
 Sardone, Nicola; Angiolini, Mauro; Belvisi, Laura;
 Maffioli, Sonia; Manzoni, Leonardo; Scolastico, Carlo
 CORPORATE SOURCE: Pharmaceutical Chemistry Department, University of
 Pavia, Pavia, 27100, Italy
 SOURCE: Tetrahedron (1998), 54(20), 5325-5336
 CODEN: TETRAB; ISSN: 0040-4020
 PUBLISHER: Elsevier Science Ltd.
 DOCUMENT TYPE: Journal
 LANGUAGE: English
 OTHER SOURCE(S): CASREACT 129:41402
 GI



AB A flexible stereoselective synthetic scheme was developed to prepare 6,5-fused bicyclic lactams I (R = Ph, 1-naphthyl, 2-naphthyl), that mol. mechanics calcns. revealed to have a potential as reverse-turn mimetics. The convergence of the synthetic sequence was achieved by attachment of a properly substituted malonate unit $\text{RCH}_2\text{CH}(\text{CO}_2\text{H})\text{CO}_2\text{Me}$ to (2S)-cis-5-(2-hydroxyethyl)proline tert-Bu ester. Stereoselective intramol. alkylation of the malonate afforded the 6-membered lactam fused to the 2-carbalkoxy pyrrolidine nucleus. X-ray crystallog. of advanced synthetic derivative II allowed the unequivocal assignment of the configuration at the newly created quaternary stereocenter as R.

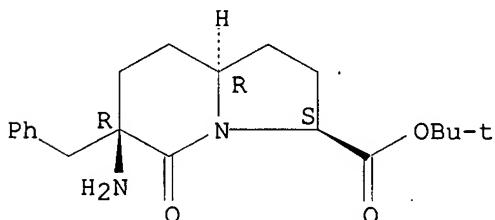
IT 208455-27-0P
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(stereoselective synthesis of bicyclic reverse-turn peptidomimetics)

RN 208455-27-0 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-aminoctahydro-5-oxo-6-(phenylmethyl)-, 1,1-dimethylethyl ester, (3S,6R,8aR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

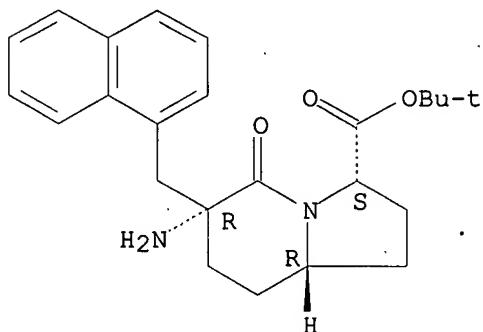


IT 208455-28-1P 208455-29-2P
 RL: SPN (Synthetic preparation); PREP (Preparation)
 (stereoselective synthesis of bicyclic reverse-turn peptidomimetics)

RN 208455-28-1 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-aminoctahydro-6-(1-naphthalenylmethyl)-5-oxo-, 1,1-dimethylethyl ester, (3S,6R,8aR)- (9CI) (CA INDEX NAME)

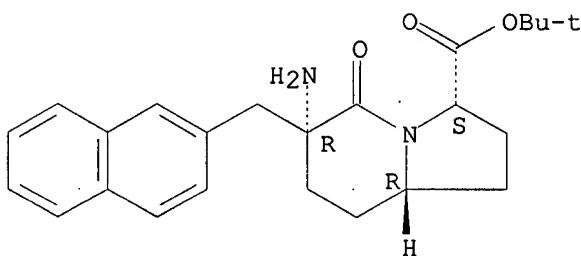
Absolute stereochemistry. Rotation (-).



RN 208455-29-2 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-aminoctahydro-6-(2-naphthalenylmethyl)-5-oxo-, 1,1-dimethylethyl ester, (3S,6R,8aR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



REFERENCE COUNT: 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 7 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1998:112235 CAPLUS

DOCUMENT NUMBER: 128:192934

TITLE: Preparation of peptide β -sheet mimetics as protease and kinase inhibitors and as inhibitors of transcription factors

INVENTOR(S): Kahn, Michael; Qabar, Maher Nicola; McMillan, Michael Kim; Ogbu, Cyprian Okwara; Eguchi, Masakatsu; Kim, Hwa-ok; Boatman, Patrick Douglas, Jr.; Urban, Jan; et al.

PATENT ASSIGNEE(S): Molecumetics Ltd., USA

SOURCE: PCT Int. Appl., 250 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

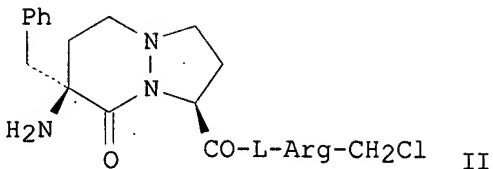
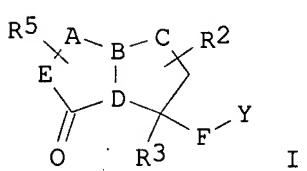
FAMILY ACC. NUM. COUNT: 3

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9805333	A1	19980212	WO 1997-US13622	19970805 <--
W: AL, AM, AT, AU, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, HU, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, AZ				

RW: GH, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, DE, DK, ES, FI, FR,
 GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA,
 GN, ML, MR, NE, SN, TD, TG
 CA 2262900 A1 19980212 CA 1997-2262900 19970805 <--
 AU 9739058 A 19980225 AU 1997-39058 19970805 <--
 AU 732174 B2 20010412
 EP 915700 A1 19990519 EP 1997-936371 19970805 <--
 EP 915700 B1 20060322
 R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
 IE, FI
 NZ 334227 A 20001027 NZ 1997-334227 19970805 <--
 JP 2001524931 T 20011204 JP 1998-508118 19970805 <--
 AT 320811 T 20060415 AT 1997-936371 19970805
 EP 1661566 A2 20060531 EP 2006-130 19970805
 R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
 IE, SI, LT, LV, FI, RO, AL
 ES 2262184 T3 20061116 ES 1997-936371 19970805
 US 6245764 B1 20010612 US 1998-9665 19980120 <--
 US 6117896 A 20000912 US 1998-22934 19980212 <--
 NO 9900522 A 19990330 NO 1999-522 19990204 <--
 KR 2000029838 A 20000525 KR 1999-700994 19990205 <--
 US 6372744 B1 20020416 US 2000-501052 20000209 <--
 US 6699869 B1 20040302 US 2000-561107 20000428
 US 2003027819 A1 20030206 US 2001-960864 20010921
 US 2004230035 A1 20041118 US 2003-745471 20031222
 US 7125872 B2 20061024
 US 2006293372 A1 20061228 US 2004-774043 20040205
 US 2006276408 A1 20061207 US 2006-448412 20060607
 PRIORITY APPLN. INFO.:
 US 1996-692420 A 19960805
 US 1996-725073 A 19961002
 US 1997-797915 A 19970210
 US 1997-47067P P 19970519
 US 1995-410518 B2 19950324
 US 1995-549006 B2 19951027
 US 1996-624690 B2 19960325
 EP 1997-936371 A3 19970805
 WO 1997-US13622 W 19970805
 US 1998-9665 A3 19980120
 US 1998-22934 A3 19980212
 US 2000-501052 A1 20000209
 US 2000-561107 A1 20000428
 US 2001-960864 B1 20010921
 US 2003-745471 A1 20031222

OTHER SOURCE(S): MARPAT 128:192934
GI



AB β -Sheet mimetics I [A = CO, (CH₂)₀₋₄, CO(CH₂)₁₋₃, (CH₂)₁₋₂CO, (CH₂)₁₋₂S; B = N, CH; C = CO, CO(CH₂)₁₋₃, (CH₂)₀₋₃, O, S, O(CH₂)₁₋₂, S(CH₂)₁₋₂; D = N, CR₄; E = CR₁NHZ, NZ, CR₁Z; F = bond, CO; R₁, R₂, R₄, R₅ = independently amino acid side chain or derivative thereof; R₂ = amino acid side chain or derivative thereof, or taken with C forms a fused substituted or unsubstituted homocyclic or heterocyclic ring; R₃ = amino acid side chain or derivative thereof, or taken with C forms a bridging moiety]

(CH₂)₁₋₂, O, S; Y, Z represent the remainder of the mol., with the proviso that any two adjacent CH groups of the bicyclic ring may form a double bond] and methods relating to the same are disclosed. The β -sheet mimetics have utility as protease and kinase inhibitors, as well as inhibitors of transcription factors. Methods of the invention include administration of a β -sheet mimetic, or use of the same for the manufacture of a medicament for treatment of a variety of conditions associated with the targeted protease, kinase and/or transcription factor. Thus, bicyclic peptide mimic II was prepared in several steps from phenylalanine Me ester, Et acrylate, and a protected arginine chloromethyl ketone derivative II was tested for inhibitory activity against a variety of serine proteases, and showed IC₅₀ = 1.2 nM against thrombin in an in vitro assay.

IT 203453-43-4P 203453-44-5P 203455-59-8P

203455-60-1P

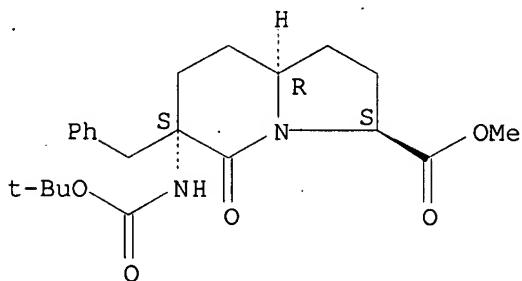
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of peptide β -sheet mimetics as protease and kinase inhibitors and as inhibitors of transcription factors)

RN 203453-43-4 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, methyl ester, [3S-(3 α ; 6 β , 8 $\alpha\beta$)]- (9CI). (CA INDEX NAME)

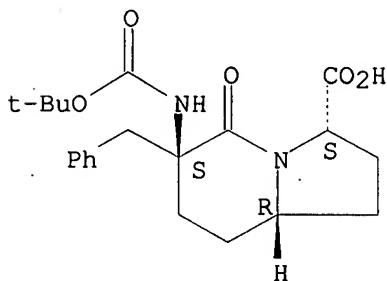
Absolute stereochemistry.



RN 203453-44-5 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, (3S, 6S, 8aR)- (9CI) (CA INDEX NAME)

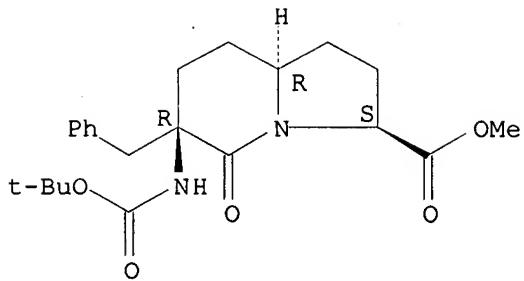
Absolute stereochemistry.



RN 203455-59-8 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, methyl ester, [3S-(3 α , 6 α , 8 $\alpha\beta$)]- (9CI) (CA INDEX NAME)

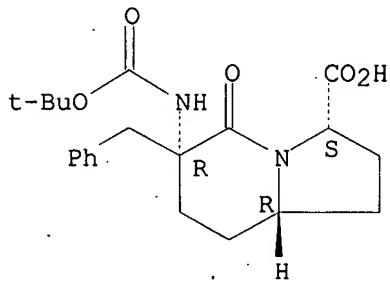
Absolute stereochemistry.



RN 203455-60-1 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, (3S,6R,8aR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 8 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1997:633831 CAPLUS

DOCUMENT NUMBER: 127:293616

TITLE: Design and synthesis of conformationally constrained arginal thrombin inhibitors

AUTHOR(S): Salimbeni, Aldo; Paleari, Fabio; Canevotti, Renato; Crisculoi, Marco; Lippi, Annalisa; Angiolini, Mauro; Belvisi, Laura; Scolastico, Carlo; Colombo, Lino
CORPORATE SOURCE: Lusofarmaco, Medicinal Chemistry Dep., Univ. Milano, Milan, 20132, Italy

SOURCE: Bioorganic & Medicinal Chemistry Letters (1997), 7(17), 2205-2210

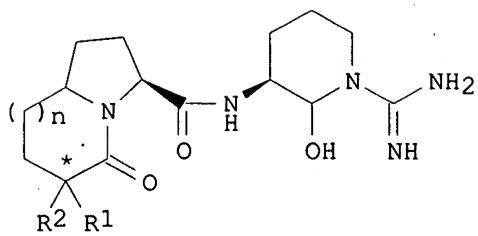
CODEN: BMCLE8; ISSN: 0960-894X

PUBLISHER: Elsevier

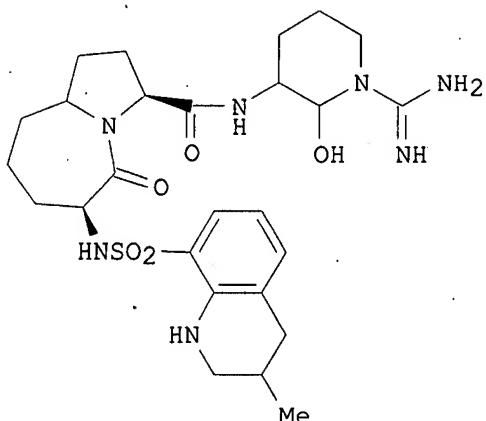
DOCUMENT TYPE: Journal

LANGUAGE: English

GI



I



II

AB A series of conformationally constrained arginal thrombin inhibitors, e.g. I ($R_1 = PhCH_2, NHAc, NHSO_2CH_2Ph$; $R_2 = H, CH_2Ph$; $n = 1, 2$; stereo at * center is RS, S, or R) was prepared starting from 5,6 or 5,7 bicyclic lactam structures, that an indirect approach of x-ray structure-based design indicated as D-Phe-Pro dipeptide mimetics. The tetrahydroquinolyl sulfonamido derivs. II (LR-D/009) displayed the best inhibitory potency ($IC_{50} = 0.018 \mu\text{m}$), with good selectivity over plasmin and trypsin.

IT 196937-03-8 196937-04-9 196937-05-0

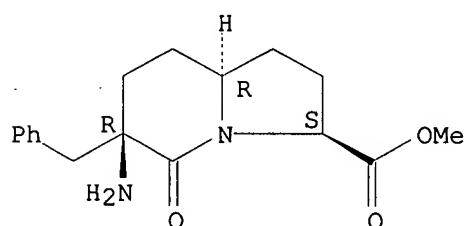
RL: PRP (Properties)

(design and synthesis of conformationally constrained arginal thrombin inhibitors)

RN 196937-03-8 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-aminooctahydro-5-oxo-6-(phenylmethyl)-, methyl ester, [3S-(3 α ,6 α ,8 β)]- (9CI) (CA INDEX NAME)

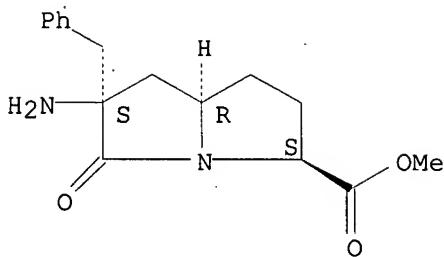
Absolute stereochemistry.



RN 196937-04-9 CAPLUS

CN 1H-Pyrrolizine-3-carboxylic acid, 6-aminohexahydro-5-oxo-6-(phenylmethyl)-, methyl ester, [3S-(3 α ,6 α ,7 β)]- (9CI) (CA INDEX NAME)

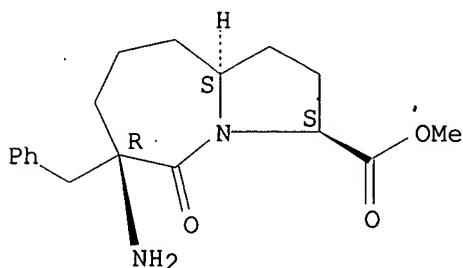
Absolute stereochemistry.



RN 196937-05-0 CAPLUS

CN 1H-Pyrrolo[1,2-a]azepine-3-carboxylic acid, 6-amino-5-oxo-6-(phenylmethyl)-, methyl ester, (3S,6R,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



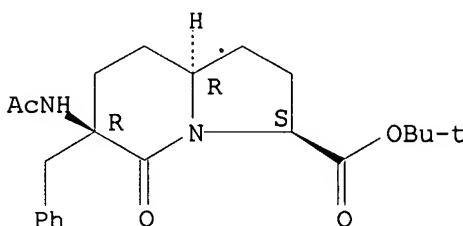
IT 162284-68-6P 188126-79-6P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(design and synthesis of conformationally constrained arginal thrombin inhibitors)

RN 162284-68-6 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-(acetylamino)octahydro-5-oxo-6-(phenylmethyl)-, 1,1-dimethylethyl ester, [3S-(3α,6α,8aβ)]- (9CI) (CA INDEX NAME)

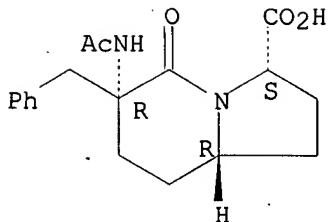
Absolute stereochemistry.



RN 188126-79-6 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-(acetylamino)octahydro-5-oxo-6-(phenylmethyl)-, [3S-(3α,6α,8aβ)]- (9CI) (CA INDEX NAME)

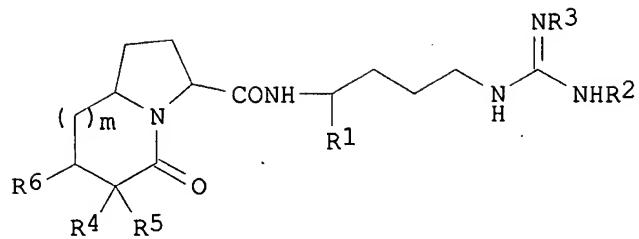
Absolute stereochemistry.



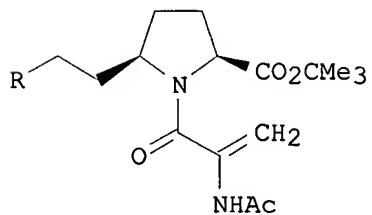
REFERENCE COUNT: 18 THERE ARE 18 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 9 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN
 ACCESSION NUMBER: 1997:231088 CAPLUS
 DOCUMENT NUMBER: 126:212450
 TITLE: Preparation of arginine-containing bicyclic lactam derivatives as thrombin inhibitors
 INVENTOR(S): Salimbeni, Aldo; Paleari, Fabio; Scolastico, Carlo; Criscuoli, Marco
 PATENT ASSIGNEE(S): A. Menarini Industrie Farmaceutiche Riunite S.R.L., Italy; Salimbeni, Aldo; Paleari, Fabio; Scolastico, Carlo; Criscuoli, Marco
 SOURCE: PCT Int. Appl., 41 pp.
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

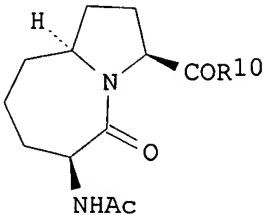
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9705160	A1	19970213	WO 1996-EP3167	19960718 <--
W: AL, AM, AT, AU, AZ, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, HU, IL, IS, JP, KE, KG, KP, KR, KZ, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE				
RW: KE, LS, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM				
AU 9667342	A	19970226	AU 1996-67342	19960718 <--
PRIORITY APPLN. INFO.:			IT 1995-MI1688	A 19950801
			WO 1996-EP3167	W 19960718
OTHER SOURCE(S): GI	MARPAT	126:212450		



I



II



III

AB Bicyclic lactams containing an arginine residue, i.e. I [$m = 0-3$; R1 = CHO, CH₂OH, CO₂H, B(OH)₂; R2, R3 = independently H, CO₂R₇, C₁₋₄ alkyl, CH₂Ph, NO₂; R4, R5 = independently H, NR₈R₉, straight or branched C₁₋₇ alkyl, C₃₋₇ cycloalkyl, or an arylalkyl or heteroarylalkyl group optionally substituted by one or more groups such as halo, OMe, CF₃, straight or branched C₁₋₇ alkyl; R6 = H, straight or branched C₁₋₇ alkyl, C₃₋₇ cycloalkyl, or an aryl, heteroaryl, arylalkyl or heteroarylalkyl group optionally substituted by one or more groups such as halo, OMe, CF₃, straight or branched C₁₋₇ alkyl; R7 = C₁₋₄ alkyl, CH₂Ph; R8, R9 = independently H, straight or branched C₁₋₇ alkyl, W-Q; W = CO, SO₂; Q = Ph, CH₂Ph, quinolyl, naphththylmethyl, tetrahydroquinolyl, optionally substituted by one or more groups such as halo, straight or branched C₁₋₇ alkyl, OMe, CF₃], which can be of use in therapy as thrombin inhibitors, are disclosed. Thus, amidation of (2S,5R)-2-tert-butoxycarbonyl-5-(2-hydroxyethyl)pyrrolidine with 2-acetylaminoacrylic acid gave 80% amide II (R = OH). Iodination of alc. II (R = OH) via its mesylate, followed by reductive radical cyclization in the presence of Bu₃SnH gave octahydropyrrolo[1,2-a]azepin-5-one III (R₁₀ = OCMe₃). Deesterification of III (R₁₀ = OCMe₃) with CF₃CO₂H, followed by coupling with N_ω-benzyloxycarbonyl-L-arginine lactam, hydride reduction, and catalytic deprotection gave arginine aldehyde derivative III (R₁₀ = L-Arg-H).

IT 162284-68-6P 188126-64-9P 188126-66-1P
188126-68-3P 188126-74-1P 188126-79-6P

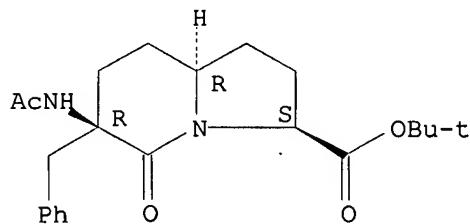
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of arginine-containing bicyclic lactam derivs. as thrombin inhibitors)

RN 162284-68-6 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-(acetylamino)octahydro-5-oxo-6-(phenylmethyl)-, 1,1-dimethylethyl ester, [3S-(3 α ,6 α ,8a β)]- (9CI) (CA INDEX NAME)

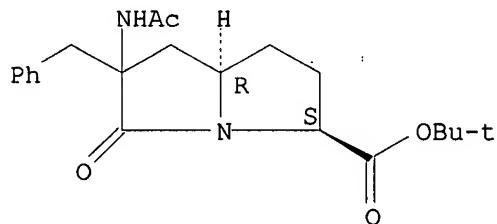
Absolute stereochemistry.



RN 188126-64-9 CAPLUS

CN 1H-Pyrrolizine-3-carboxylic acid, 6-(acetylamino)hexahydro-5-oxo-6-(phenylmethyl)-, 1,1-dimethylethyl ester, (3S,7aR)- (9CI) (CA INDEX NAME)

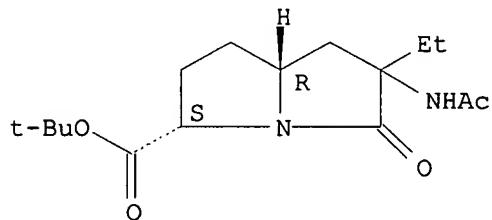
Absolute stereochemistry.



RN 188126-66-1 CAPLUS

CN 1H-Pyrrolizine-3-carboxylic acid, 6-(acetylamino)-6-ethylhexahydro-5-oxo-, 1,1-dimethylethyl ester, (3S,7aR)- (9CI) (CA INDEX NAME)

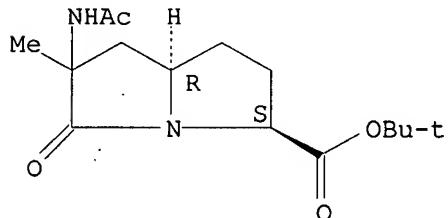
Absolute stereochemistry.



RN 188126-68-3 CAPLUS

CN 1H-Pyrrolizine-3-carboxylic acid, 6-(acetylamino)hexahydro-6-methyl-5-oxo-, 1,1-dimethylethyl ester, (3S,7aR)- (9CI) (CA INDEX NAME)

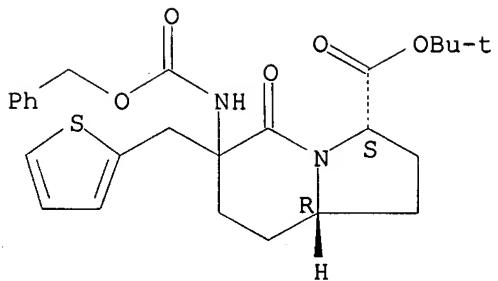
Absolute stereochemistry.



RN 188126-74-1 CAPLUS

CN 3-Indolizinecarboxylic acid, octahydro-5-oxo-6-[(phenylmethoxy)carbonyl]amino]-6-(2-thienylmethyl)-, 1,1-dimethylethyl ester, (3S,8aR)-[partial]- (9CI) (CA INDEX NAME)

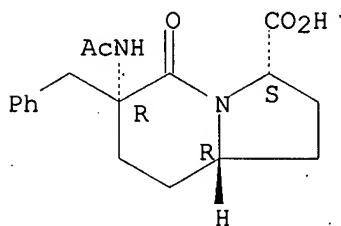
Absolute stereochemistry.



RN 188126-79-6 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-(acetylaminoo)octahydro-5-oxo-6-(phenylmethyl)-, [3S-(3 α ,6 α ,8a β)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



L4 ANSWER 10 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1996:731812 CAPLUS

DOCUMENT NUMBER: 126:8708

TITLE: Preparation of beta-sheet mimetics of peptides or proteins as protease inhibitors

INVENTOR(S): Kahn, Michael

PATENT ASSIGNEE(S): Molecumetics Ltd., USA

SOURCE: PCT Int. Appl., 141 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

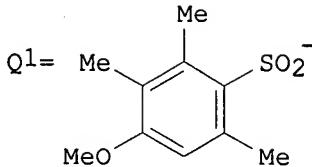
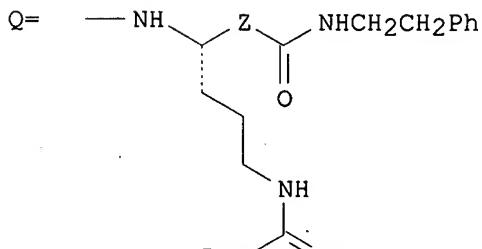
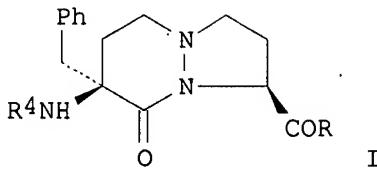
FAMILY ACC. NUM. COUNT: 3

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9630396	A1	19961003	WO 1996-US4115	19960325 <--
W: AL, AM, AT, AU, AZ, BB, BG, BR, BY, CA, CH, CN, CZ, DE, DK, EE, ES, FI, GB, GE, HU, IS, JP, KE, KG, KP, KR, KZ, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI				
RW: KE, LS, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN				
AU 9653729	A	19961016	AU 1996-53729	19960325 <--
AU 713530	B2	19991202		
EP 815123	A1	19980107	EP 1996-910566	19960325 <--
EP 815123	B1	20011004		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, FI				
JP 10508035	T	19980804	JP 1996-529594	19960325 <--
JP 2000319295	A	20001121	JP 2000-79170	19960325 <--
AT 206433	T	20011015	AT 1996-910566	19960325 <--
US 6020331	A	20000201	US 1998-9386	19980120 <--
US 6245764	B1	20010612	US 1998-9665	19980120 <--
US 6586426	B1	20030701	US 1999-443055	19991118

US 6699869	B1	20040302	US 2000-561107	20000428
US 2003191109	A1	20031009	US 2001-8770	20011025
US 2004230035	A1	20041118	US 2003-745471	20031222
US 7125872	B2	20061024		
JP 2004131511	A	20040430	JP 2004-17703	20040126
US 2006276408	A1	20061207	US 2006-448412	20060607
PRIORITY APPLN. INFO.:				
			US 1995-410518	A 19950324
			US 1995-549006	A 19951027
			JP 1996-529594	A3 19960325
			JP 2000-79170	A3 19960325
			US 1996-624690	B2 19960325
			US 1996-624695	B1 19960325
			WO 1996-US4115	W 19960325
			US 1996-725073	B1 19961002
			US 1998-4968	B1 19980109
			US 1998-9386	A3 19980120
			US 1998-9665	A3 19980120
			US 2000-561107	A1 20000428
			US 2003-745471	A1 20031222

OTHER SOURCE(S): MARPAT 126:8708
GI



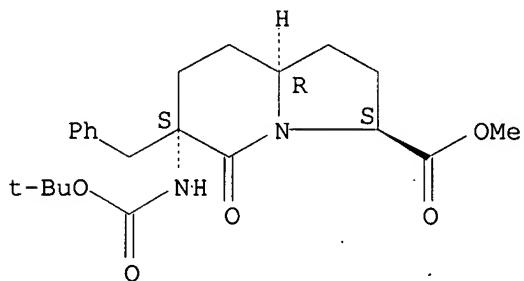
AB There are disclosed β -sheet mimetics [I; R1 - R3 = amino acid side chain moiety or its derivative; A = CO, $(\text{CH}_2)_1-4$, $(\text{CH}_2)_1-2-\text{O}$, $(\text{CH}_2)_1-2-\text{S}$; B = N, CH; C = CO, $(\text{CH}_2)_1-3$, O, S, $\text{O}(\text{CH}_2)_1-2$, $\text{S}(\text{CH}_2)_1-2$; Y, Z = the remainder of the mol.; or any 2 adjacent CH groups of the bicyclic ring may form a double bond; with the provisos that (1) R1 = an amino acid side chain moiety or derivative thereof other than H, (2) when R1 = CH_2Ph , R2 = R3 = H, A = CH_2CH_2 , and B = CH, and then C \neq CH2, (3) when R1 = me, R2 and R3 = H, A = CH_2O , B = CH, and then C \neq CH2, and (4) when R1 = CH_2Ph , R2 = R3 = H, A = CH2, B = CH, and then C \neq S] and methods relating to the same for imparting or stabilizing the β -sheet structure of a peptide, protein or mol. In one aspect, the β -sheet mimetics are covalently attached at the end or within the length of the peptide or protein. The β -sheet mimetics have utility as protease inhibitors generally, including activity as serine protease inhibitors such as thrombin, elastase, and Factor X. Thus, diazabicyclo[4.3.0]nonane derivative (II; R = OH, R4 = Boc) (preparation given) was condensed with amino alc. (H-Q; R5 = Q1, Z = CHOH) using 1-ethyl-3-(3-dimethylaminopropyl)carbodiimide hydrochloride, HOEt hydrate, and $(\text{Me}_2\text{CH})_2\text{NEt}$ in THF to give amide alc. II (R = Q, R5 = Q1, Z = CHOH, R4 = Boc), which was oxidized by Dess-Martin periodinane in CH_2Cl_2 and deprotected with 95% aqueous $\text{CF}_3\text{CO}_2\text{H}$ and thioanisole to give the β -sheet mimetic II (R = Q, R5 = H, Z = CO, R4 = H). The latter compound in vitro inhibited various serine proteases such as.

thrombin, factor VII, factor X, and trypsin, e.g. with K_i of 7.10 + 10-11 M for thrombin.

IT 183442-92-4P 183442-93-5P 183442-97-9P
 183624-03-5P
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
 (preparation of beta-sheet mimetics of peptides or proteins as protease inhibitors)

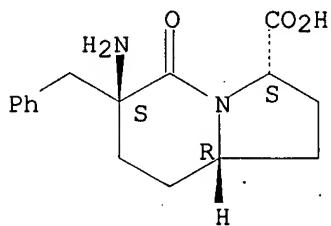
RN 183442-92-4 CAPLUS
 CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, methyl ester, (3R,6R,8aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



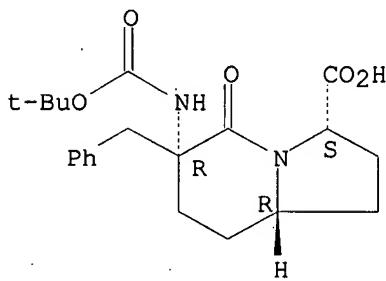
RN 183442-93-5 CAPLUS
 CN 3-Indolizinecarboxylic acid, 6-amino-octahydro-5-oxo-6-(phenylmethyl)-, (3R,6R,8aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 183442-97-9 CAPLUS
 CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, (3α,6α,8aβ)- (9CI) (CA INDEX NAME)

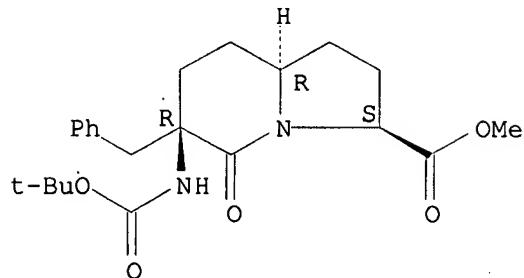
Relative stereochemistry.



RN 183624-03-5 CAPLUS
 CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, methyl ester, (3R,6S,8aS)-rel- (9CI) (CA INDEX NAME)

INDEX NAME)

Relative stereochemistry.



L4 ANSWER 11 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1996:731810 CAPLUS

DOCUMENT NUMBER: 126:8707

TITLE: Preparation of beta-sheet mimetics of peptides or proteins as inhibitors of biologically active peptides or proteins

INVENTOR(S): Kahn, Michael

PATENT ASSIGNEE(S): Molecumetics Ltd., USA

SOURCE: PCT Int. Appl., 158 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 3

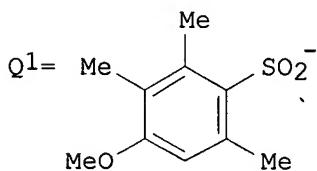
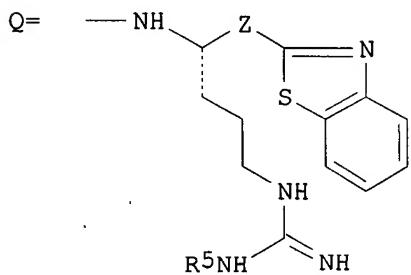
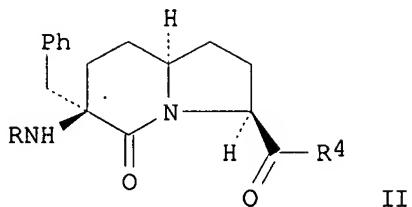
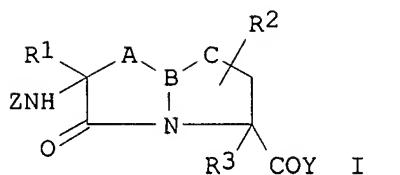
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9630035	A1	19961003	WO 1996-US4044	19960325 <--
W: AL, AM, AT, AU, AZ, BB, BG, BR, BY, CA, CH, CN, CZ, DE, DK, EE, ES, FI, GB, GE, HU, IS, JP, KE, KG, KP, KR, KZ, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI				
RW: KE, LS, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN				
CA 2215695	A1	19961003	CA 1996-2215695	19960325 <--
CA 2215695	C	20030916		
CA 2215720	A1	19961003	CA 1996-2215720	19960325 <--
AU 9653714	A	19961016	AU 1996-53714	19960325 <--
AU 712581	B2	19991111		
EP 817642	A1	19980114	EP 1996-910547	19960325 <--
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, FI				
JP 10508034	T	19980804	JP 1996-529567	19960325 <--
JP 2000319295	A	20001121	JP 2000-79170	19960325 <--
ES 2161354	T3	20011201	ES 1996-910566	19960325 <--
US 6020331	A	20000201	US 1998-9386	19980120 <--
US 6245764	B1	20010612	US 1998-9665	19980120 <--
US 6586426	B1	20030701	US 1999-443055	19991118
US 6699869	B1	20040302	US 2000-561107	20000428
US 2003191109	A1	20031009	US 2001-8770	20011025
US 2004230035	A1	20041118	US 2003-745471	20031222
US 7125872	B2	20061024		
JP 2004131511	A	20040430	JP 2004-17703	20040126
US 2006276408	A1	20061207	US 2006-448412	20060607
PRIORITY APPLN. INFO.:			US 1995-410518	A 19950324
			US 1995-549006	A 19951027
			JP 1996-529594	A3 19960325
			JP 2000-79170	A3 19960325

US 1996-624690	B2 19960325
US 1996-624695	B1 19960325
WO 1996-US4044	W 19960325
US 1996-725073	B1 19961002
US 1998-4968	B1 19980109
US 1998-9386	A3 19980120
US 1998-9665	A3 19980120
US 2000-561107	A1 20000428
US 2003-745471	A1 20031222

OTHER SOURCE(S) :
GI

MARPAT 126:8707



AB There are disclosed β -sheet mimetics [I; R1 - R3 = amino acid side chain moiety or its derivative; A = CO, (CH₂)₁₋₄, (CH₂)_{1-2-O}, (CH₂)_{1-2-S}; B = N, CH; C = CO, (CH₂)₁₋₃, O, S, O(CH₂)₁₋₂, S(CH₂)₁₋₂; Y, Z = the remainder of the mol.; or any 2 adjacent CH groups of the bicyclic ring may form a double bond] and methods relating to the same for imparting or stabilizing the β -sheet structure of a peptide, protein or mol. In one aspect, the β -sheet mimetics are covalently attached at the end or within the length of the peptide or protein. The β -sheet mimetics have utility as inhibitors of one or more of proteases, kinases, CAAX motif (Ras prenylation of the Cys within its C-terminal CAAX sequence by farnesyl transferase, wherein "A" is defined as an amino acid with a hydrophobic side chain and "X" is another amino acid), peptides binding to SH2 domains, and MHC-I and/or MHC-II (major histocompatibility complex class I and class II) presentation of peptides to T cell receptors in warm-blooded animals. Thus, azabicyclo[4.3.0]nonane derivative (II; R = Boc, R4 = OH) (preparation given) was condensed with benzothiazolylarginol derivative (H-Q.CF₃CO₂H; R5 = Q1, Z = CHOH) using 1-ethyl-3-(3-dimethylaminopropyl)carbodiimide hydrochloride, HOEt, and (Me₂CH)₂NEt in THF to give arginol derivative II (R = Boc, R4 = Q, R5 = Q1 Z = CHOH), which was oxidized by Dess-Martin periodinane in CH₂C₁₂ to arginine derivative II (R = Boc, R4 = Q, R5 = Q1 Z = CO) and deprotected 95% aqueous CF₃CO₂H and thioanisole at room temperature for 20 h to give, after HPLC purification, the β -sheet mimetic II (R = H, R4 = Q, R5 = H, Z = CO). The latter compound in vitro inhibited various serine proteases such as thrombin, factor VII, factor X, factor XI, urokinase, thrombin-thrombomodulin complex, activated protein C, plasmin, tissue plasminogen activator, trypsin, and tryptase, e.g. with Ki of 8.50 + 10⁻¹¹ M for thrombin.

IT 183442-92-4P 183442-93-5P 183442-97-9P
183624-03-5P

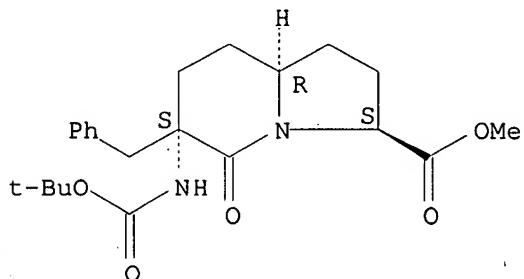
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of beta-sheet mimetics of peptides or proteins as inhibitors of biol. active peptides or proteins)

RN 183442-92-4 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, methyl ester, (3R,6R,8aS)-rel- (9CI) (CA INDEX NAME)

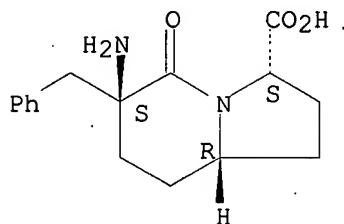
Relative stereochemistry.



RN 183442-93-5 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-aminooctahydro-5-oxo-6-(phenylmethyl)-, (3R,6R,8aS)-rel- (9CI) (CA INDEX NAME)

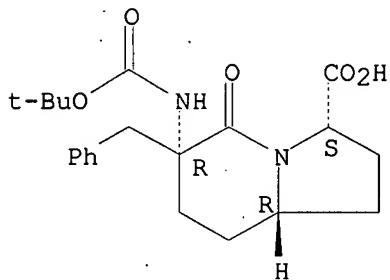
Relative stereochemistry.



RN 183442-97-9 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, (3α,6α,8aβ)- (9CI) (CA INDEX NAME)

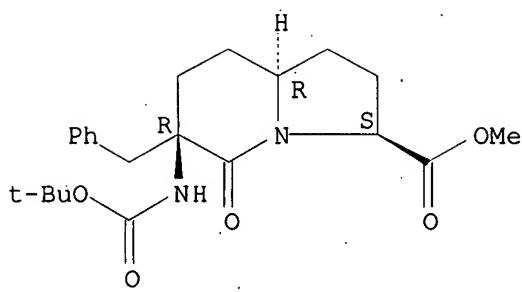
Relative stereochemistry.



RN 183624-03-5 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-[[[(1,1-dimethylethoxy)carbonyl]amino]octahydro-5-oxo-6-(phenylmethyl)-, methyl ester, (3R,6S,8aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



L4 ANSWER 12 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1996:662548 CAPLUS

DOCUMENT NUMBER: 126:19211

TITLE: Conformationally constrained dipeptides: synthesis of bicyclic lactams by stereoselective radical cyclizations

AUTHOR(S): Colombo, Lino; Di Giacomo, Marcello; Belvisi, Laura; Manzoni, Leonardo; Scolastico, Carlo; Salimbeni, Aldo

CORPORATE SOURCE: Dip. Chim. Farmaceutica, Univ. Pavia, Pavia, I-27100, Italy

SOURCE: Gazzetta Chimica Italiana (1996), 126(8), 543-554

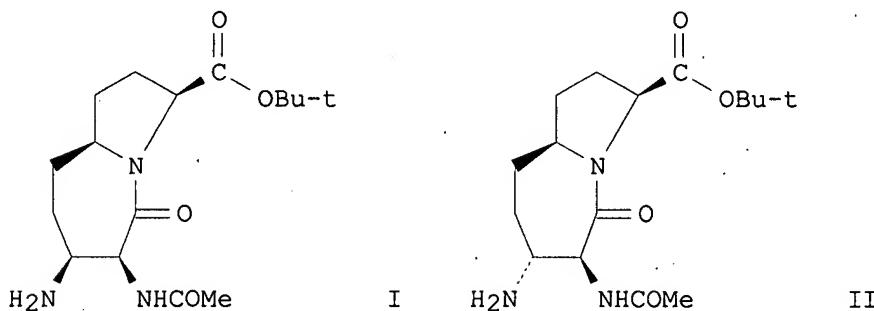
PUBLISHER: CODEN: GCITA9; ISSN: 0016-5603
Societa Chimica Italiana

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 126:19211

GI



AB Regioselective and stereoselective radical cyclization of β -substituted α -N-acetyl acrylamides was performed. This method led to the formation of 6,5- and 7,5-fused bicyclic lactams, which can be viewed as conformationally restricted dipeptide mimics. The peptide mimics I and II were prepared

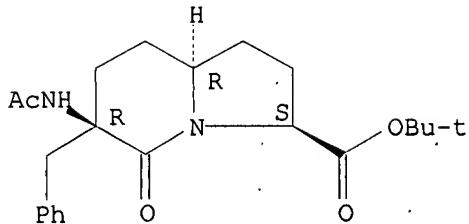
IT 162284-68-6P

RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of conformationally constrained dipeptide mimics by radical cyclization of β -substituted α -N-acetyl acrylamides)

RN 162284-68-6 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-(acetylamino)octahydro-5-oxo-6-(phenylmethyl)-, 1,1-dimethylethyl ester, [3S-(3 α ,6 α ,8 α B)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



L4 ANSWER 13 OF 13 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1995:357314 CAPLUS

DOCUMENT NUMBER: 122:240379

TITLE: Conformationally constrained dipeptides: synthesis of 7,5- and 6,5-fused bicyclic lactams by stereoselective radical cyclizations

AUTHOR(S): Colombo, Lino; Di Giacomo, Marcello; Scolastico, Carlo; Manzoni, Leonardo; Belvisi, Laura; Molteni, Valentina

CORPORATE SOURCE: Dipart. Chim. Farm., Univ. Pavia, Pavia, I-27100, Italy

SOURCE: Tetrahedron Letters (1995), 36(4), 625-8
CODEN: TELEAY; ISSN: 0040-4039

PUBLISHER: Elsevier

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 122:240379

AB A study of radical cyclizations of β -substituted α -N-acetyl acrylamide have been performed: high level of regio- and stereoselectivity was obtained.

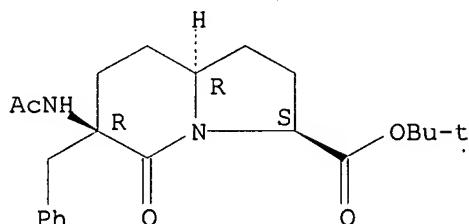
IT 162284-68-6P

RL: SPN (Synthetic preparation); PREP (Preparation)
(conformationally constrained dipeptides and synthesis of fused bicyclic lactams by stereoselective radical cyclizations)

RN 162284-68-6 CAPLUS

CN 3-Indolizinecarboxylic acid, 6-(acetylamino)octahydro-5-oxo-6-(phenylmethyl)-, 1,1-dimethylethyl ester, [3S-(3 α ,6 α ,8 $\alpha\beta$)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



=> FIL STNGUIDE

COST IN U.S. DOLLARS

SINCE FILE ENTRY	TOTAL SESSION
---------------------	------------------

FULL ESTIMATED COST

71.46 243.77

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE ENTRY	TOTAL SESSION
---------------------	------------------

CA SUBSCRIBER PRICE

-10.14 -10.14

FILE 'STNGUIDE' ENTERED AT 11:17:53 ON 02 JUL 2007

USE IS SUBJECT TO THE TERMS OF YOUR CUSTOMER AGREEMENT
COPYRIGHT (C) 2007 AMERICAN CHEMICAL SOCIETY (ACS)

FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Jun 29, 2007 (20070629/UP).

=> d his

(FILE 'HOME' ENTERED AT 11:16:18 ON 02 JUL 2007)

FILE 'REGISTRY' ENTERED AT 11:16:27 ON 02 JUL 2007
L1 STRUCTURE uploaded
L2 43 S L1 FULL

FILE 'CAPLUS' ENTERED AT 11:16:54 ON 02 JUL 2007
L3 18 S L2 FULL
L4 13 S L3 AND PY<2003

FILE 'STNGUIDE' ENTERED AT 11:17:53 ON 02 JUL 2007

=> log y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.24	244.01
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-10.14

STN INTERNATIONAL LOGOFF AT 11:20:34 ON 02 JUL 2007